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Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (currently amended) An acid addition salt of a compound of Formula I

wherein R² represents C₁₋₆ alkyl (optionally substituted and/or terminated by one or more substituents selected from -OH, halo, cyano, nitro and aryl) or aryl, wherein each aryl and aryloxy group, unless otherwise specified, is optionally substituted.

2. (currently amended) A salt according to claim 1, wherein in which the acid component of the acid addition salt is represented by formula A

wherein R^{16} represents unsubstituted C_{1-4} alkyl, C_{1-4} perfluoroalkyl or phenyl, which latter group is optionally substituted by one or more substituents selected from C_{1-6} alkyl, halo, nitro and C_{1-6} alkoxy, and R^2 is as defined above.

- 3. (currently amended) A salt according to claim 2, wherein the salt is a toluenesulfonate, benzenesulfonate, nosylate, brosylate, besylate or mesitylate salt.
- 4. (currently amended) A salt according to any previous claim 1, wherein in which the salt is in solid form.

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5. (currently amended) A salt according to any previous claim 1, wherein the salt which is [2-(9-oxa-3,7-diazabicyclo[3.3.1]non-3-yl)-ethyl]-carbamic acid *tert*-butyl ester 2,4,6-trimethylbenzenesuifonic acid.

6. (currently amended) A process for the preparation of a compound of Formula II

wherein R¹ represents a structural fragment moiety of formula Ia

in which wherein A represents CH₂ and R³ represents -OH or -N(H)R⁷;

 R^4 represents H, C_{1-6} alkyl or, together with R^3 , represents =0;

 R^5 represents phenyl or pyridyl, both of which groups are optionally substituted by one or more substituents selected from -OH, cyano, halo, nitro, C_{1-6} alkyl (optionally terminated by - $N(H)C(O)OR^{13a}$), C_{1-6} alkoxy, $-N(R^{14a})R^{14b}$, $-C(O)R^{14c}$, $-C(O)OR^{14d}$, $-C(O)N(R^{14e})R^{14f}$, - $N(R^{14g})C(O)R^{14h}$, $-N(R^{14i})C(O)N(R^{14j})R^{14k}$, $-N(R^{14m})S(O)_2R^{13b}$, $-S(O)_2R^{13c}$ and/or - $OS(O)_2R^{13d}$;

 R^7 represents H, C_{1-6} alkyl, -E-aryl, -E-Het¹, -C(O) R^{9a} , -C(O)O R^{9b} , -S(O)₂ R^{9c} , - [C(O)]_pN(R^{10a}) R^{10b} or -C(NH)NH₂;

- R^{9a} to R^{9d} independently represent, independently at each occurrence when used herein, C_{1-6} alkyl (optionally substituted and/or terminated by one or more substituents selected from halo, aryl and Het^2), aryl, Het^3 , or R^{9a} and R^{9d} independently represent H;
- R^{10a} and R^{10b} independently represent, at each occurrence when used herein, H or C_{1-6} alkyl (optionally substituted and/or terminated by one or more substituents selected from halo, aryl and Het⁴), aryl, Het⁵, or together represent C_{3-6} alkylene, optionally interrupted by an O atom;

E represents, independently at each occurrence when used herein, a direct bond or C₁₋₄ alkylene;

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B represents -Z-, -Z-N(R^{12})-, -N(R^{12})-Z-, -Z-S(O)_n- or -Z-O- (in which latter two groups, Z is attached to the carbon atom bearing R^3 and R^4);

 \overline{Z} represents a direct bond or \overline{C}_{1-4} alkylene;

R¹¹ and R¹² independently represent H or C₁₋₆ alkyl;

R^{13a} to R^{13d} independently represent C₁₋₆ alkyl;

R^{14a} and R^{14b} independently represent H, C₁₋₆ alkyl or together represent C₃₋₆ alkylene, resulting in a four- to seven-membered nitrogen-containing ring;

R^{14c} to R^{14m} independently represent H or C₁₋₆ alkyl; and

n represents 0, 1 or 2;

p represents 1 or 2;

Het¹ to Het⁵ independently represent, independently at each occurrence when used herein, five-to twelve-membered heterocyclic groups containing one or more heteroatoms selected from oxygen, nitrogen and/or sulfur, which heterocyclic groups are optionally substituted by one or more substituents selected from =O, -OH, cyano, halo, nitro, C₁₋₆ alkyl, C₁₋₆ alkoxy, aryl, aryloxy, -N(R^{15a})R^{15b}, -C(O)R^{15c}, -C(O)OR^{15d}, -C(O)N(R^{15e})R^{15f}, -N(R^{15g})C(O)R^{15h} and -N(R¹⁵ⁱ)S(O)₂R^{15j};

R^{15a} to R^{15j} independently represent C₁₋₆ alkyl, aryl or R^{15a} to R¹⁵ⁱ independently represent H; and R² represents C₁₋₆ alkyl (optionally substituted and/or terminated by one or more substituents selected from -OH, halo, cyano, nitro and aryl) or aryl, wherein each aryl and aryloxy group, unless otherwise specified, is optionally substituted.

wherein a salt of a compound of Formula I

$$\begin{array}{c|c}
 & O \\
 & O \\$$

 $\underline{\text{in which}}\ \underline{\text{wherein}}\ R^2$ is as previously defined is reacted with a compound of Formula III

wherein Y represents O or N(R⁷) and R⁴, R⁵, R⁷ and B are as hereinbefore defined,

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at a temperature in the range of 0 °C to 100 °C for example at elevated temperature (e.g. 60°C to reflux) in the presence of a water and in the presence of a base.

- 7. (currently amended) A process according to claim 6, wherein in which the salt has been previously isolated in solid form.
- 8. (currently amended) A process according to either claim 6 or claim 7 for the preparation of tert-butyl 2-{7-[(2S)-3-(4-cyanophenoxy)-2-hydroxypropyl]-9-oxa-3,7-diaza-bicyclo[3.3.1]-non-3-yl}ethylcarbamate which comprises reacting a salt of [2-(9-oxa-3,7-diaza-bicyclo[3.3.1]non-3-yl)-ethyl]-carbamic acid tert-butyl ester with 4-[(2S)-oxiranylmethoxy]benzonitrile at a temperature in the range of 0 °C to 100 °C in the presence of water and in the presence of a base.
- 9. (currently amended) A process according to any one of claims claim 6, wherein 7 or 8 in which the salt of Formula I is an isolated salt of [2-(9-oxa-3,7-diazabicyclo[3.3.1]non-3-yl)-ethyl]-carbamic acid *tert*-butyl ester is used.
- 10. (currently amended) A process according to either claim 9, wherein the salt is the 2,4,6-trimethylbenzenesulfonic acid salt.